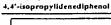




SDBS-13C NMRSDBS No. 1716CDS-12-007  $\mathrm{C_{13}\,H_{16}\,O_2}$ 

22.53 MHz 9.025 g : 0.5 ml OMSO-d<sub>4</sub>



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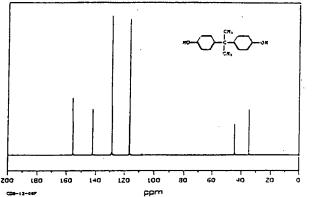
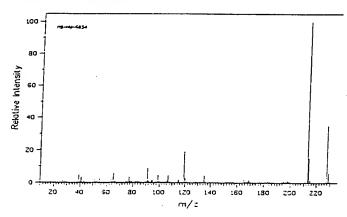


Fig. I (a)

Fig. 2 (6)



MS-MM-9834 IDBS MO. 1716 1,47-150PPGPYLIDEMEDIPHENOU CISM1602 IMags of molecular ton: 2299



Source Temperature: 190 C Sample Temperature: 140 C Direct, 75 eV

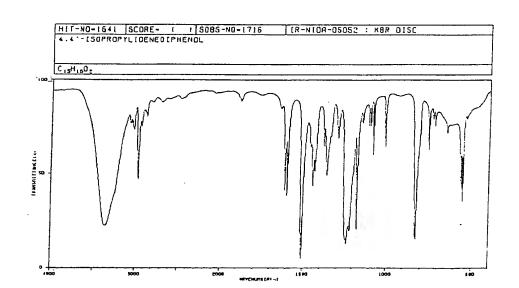


Fig. 1 (c)

### SDBS-Mass

223)

Fig 2(b)

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13

27.0 39.0 41.0 351.0 255.0 63.0 65.0 66.0 76.0 77.0 379.0 299.0 91.0 99.0 91.0 99.0 91.0 91.0 91.0

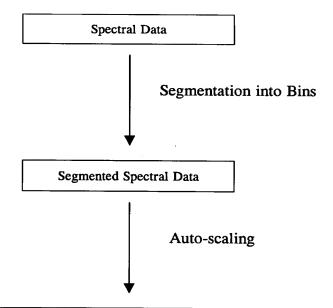
<u> </u>	31	Cm	31	Chai	31	_Cira_	21	CM	21	•
	21	2933	72	1436	49			816	66	
	74	2871		1384	62			759	50	CH,
050	74	1512	39	1363	47	1113	72	735		
	70	1600	37	1296		1102			72	но— — С— — Он
976	46	1210	4	1247		1085		650	sa	PO-C-V )-OH
966		1453	62	1239	12	1013	62	565	41	
956	62	1447	42	1221		827		663	34	ĊH-

Fig 2(c)

Bin Number	Spectral	Bin Number	Spectral	Bin Number	Spectral
	Intensity		Intensity		Intensity
51	2	183	1	1025	62
55	2	195	1	1026	42
63	2	197	1	1027	49
65	5	198	1	1032	62
66	l	212	1 .	1034	47
76	I	213	100	1041	66
77	3	214	15	1046	14
79	1	215	1	1047	12
89	I	228	35	1048	19
90	1	229	5	1053	20
91	9	580	328	1056	74
92	ı	590	223	1059	72
94	2	664	977	1060	72
95	1	677	1000	1062	57
99	4	690	330	1069	62
105	I	704	441	1088	14
106	2	834	21	1089	55
107	4	864	74	1095	60
114	ı	866	74	1097	77
15	1	868	70	1097	72
19	19	873	46	1106	
20	2	874	46	1114	68
34	i	875	62		41
35	4	877	72	1115	34
52	<del>                                     </del>	883	77		
65	2	1009	39		
69	ī	1011	37		
81	<del>  i                                   </del>	1011			
· · · · · · · · · · · · · · · · · · ·	╙・──	1 1017	4	1	

Figure 3 – a hypothetical set of spectrally derived molecular structure descriptors for bisphenol A.

FIG. 4



Auto-scaled, segmented spectral data that equalizes the importance of inherently weak spectral signals and inherently strong spectral signals.

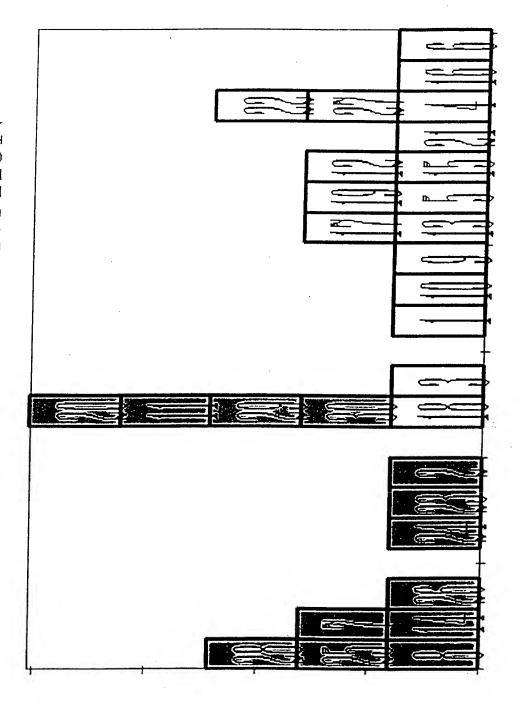
Fisher-Weighting

Fisher-weighted, auto-scaled, segmented spectral data that emphasizes those spectral characteristics that are most important for discerning the endpoint classes of the compounds in the training set.

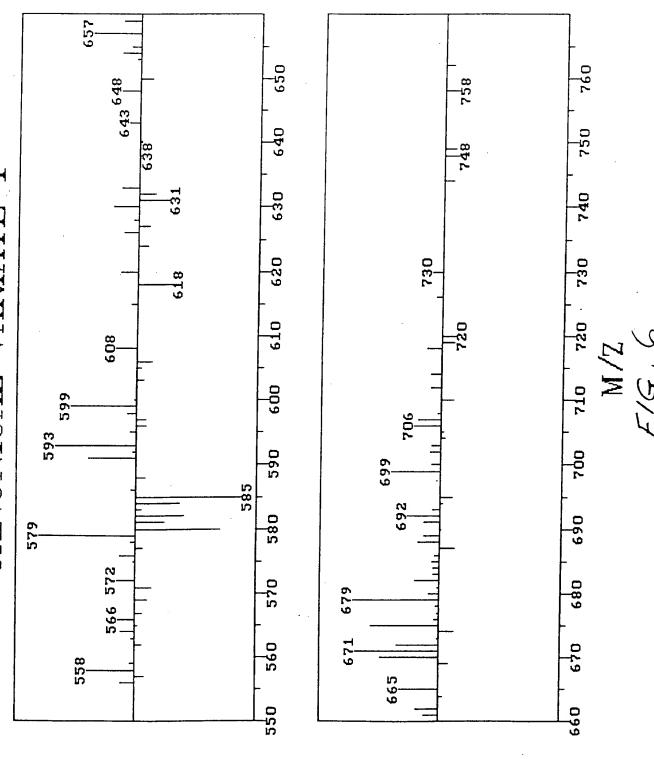
Pattern Recognition by, for example, Principal Component Analysis (PCA)

The spectral data-activity relationship (SDAR), a quantitative relationship between segmented spectral data and endpoint data. The SDAR may be visualized with a canonical variate function, that illustrates how spectral data in a particular bin shows bias toward a particular endpoint property and a discriminant function that illustrates how well the SDAR is able to discern the endpoint group of a compound from its segmented spectral data.

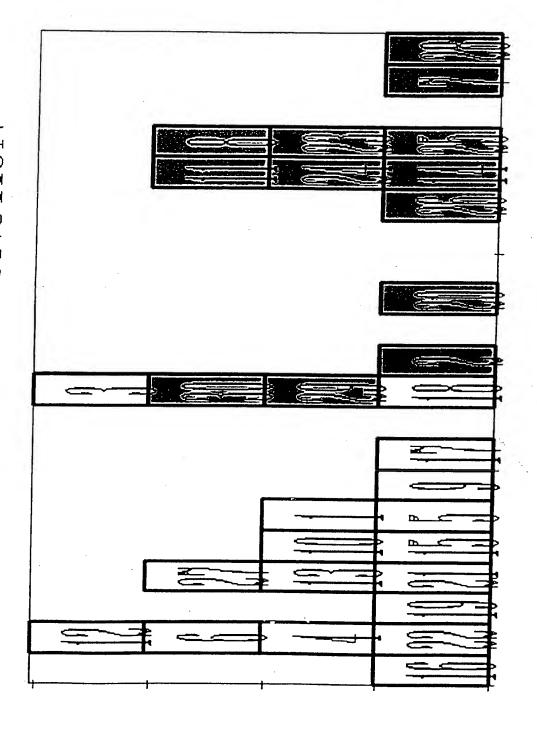
# DISCRIMINANT FUNCTION



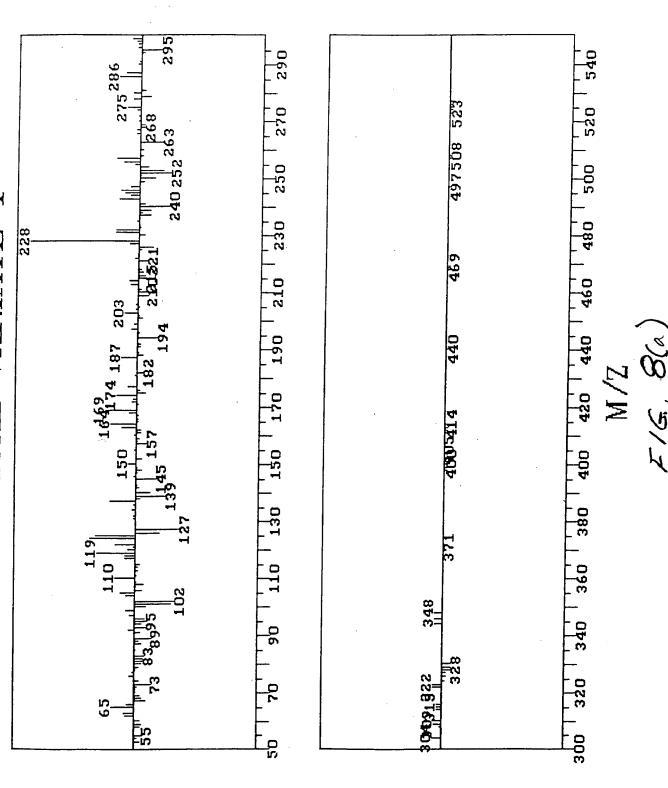
### COMPONENT 1

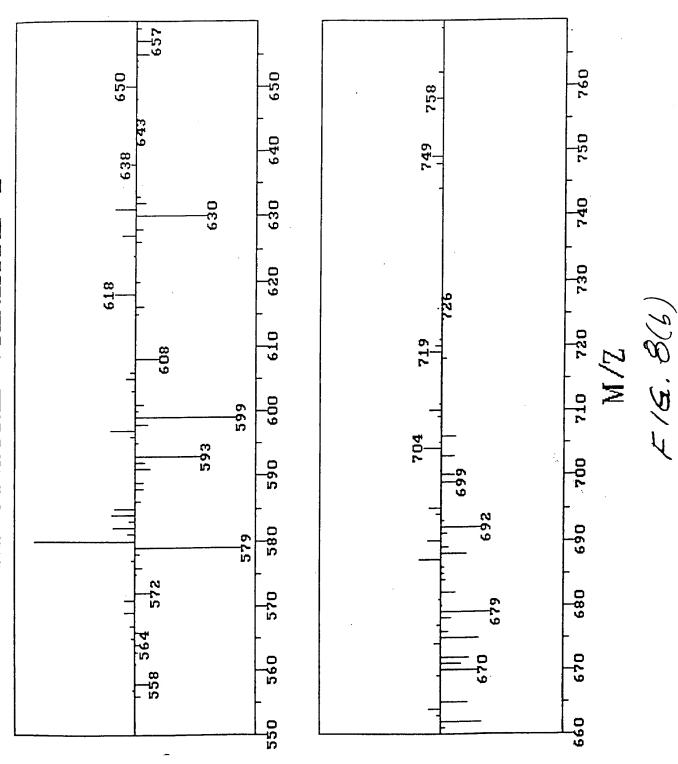


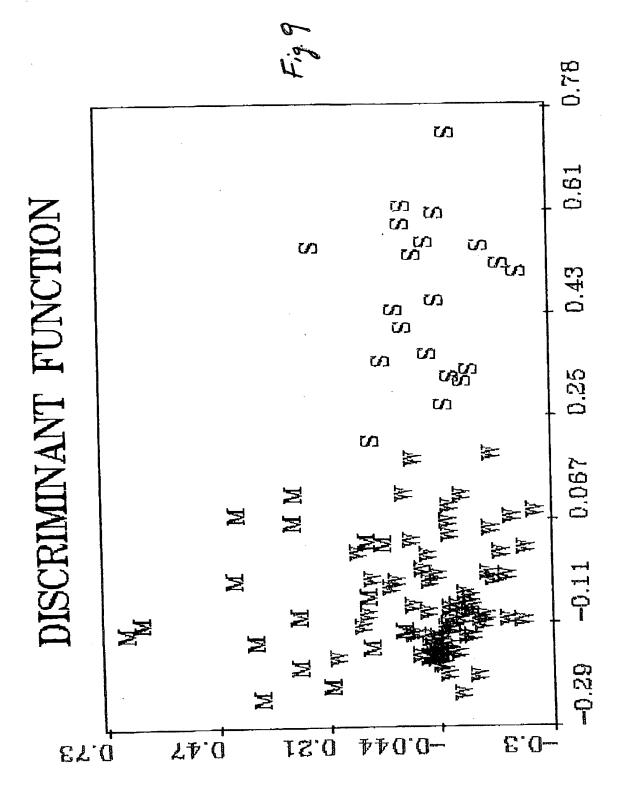
# DISCRIMINANT FUNCTION

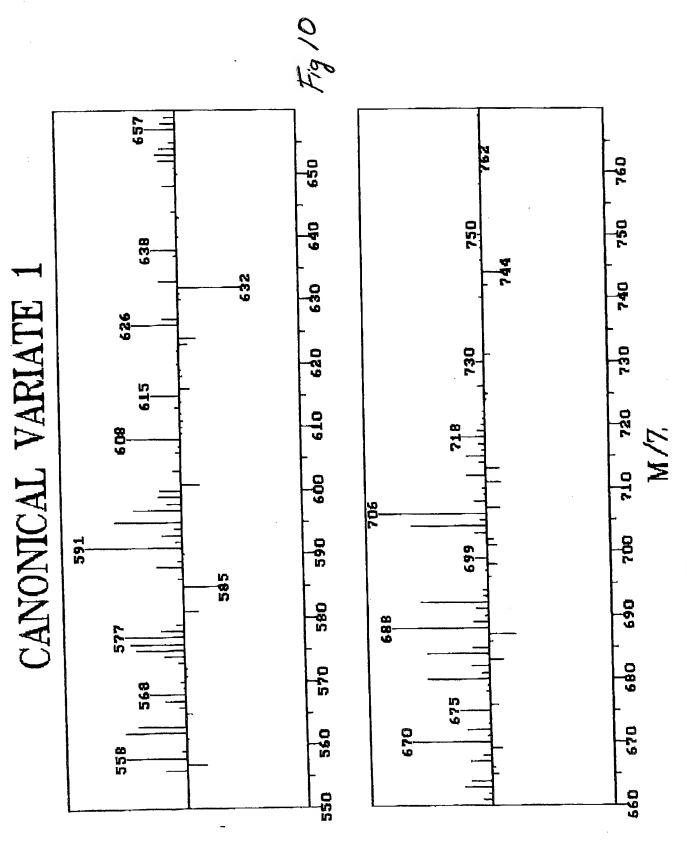


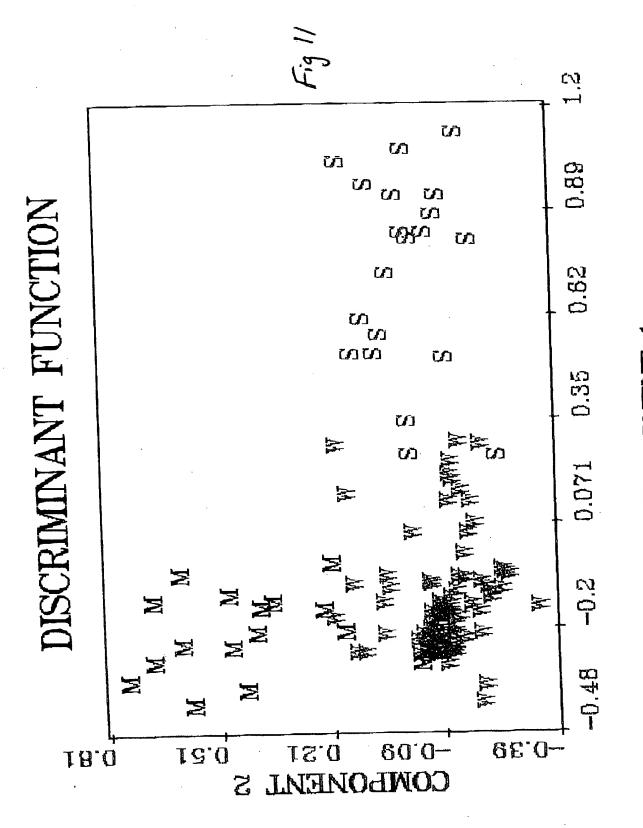
COMPONENT 1



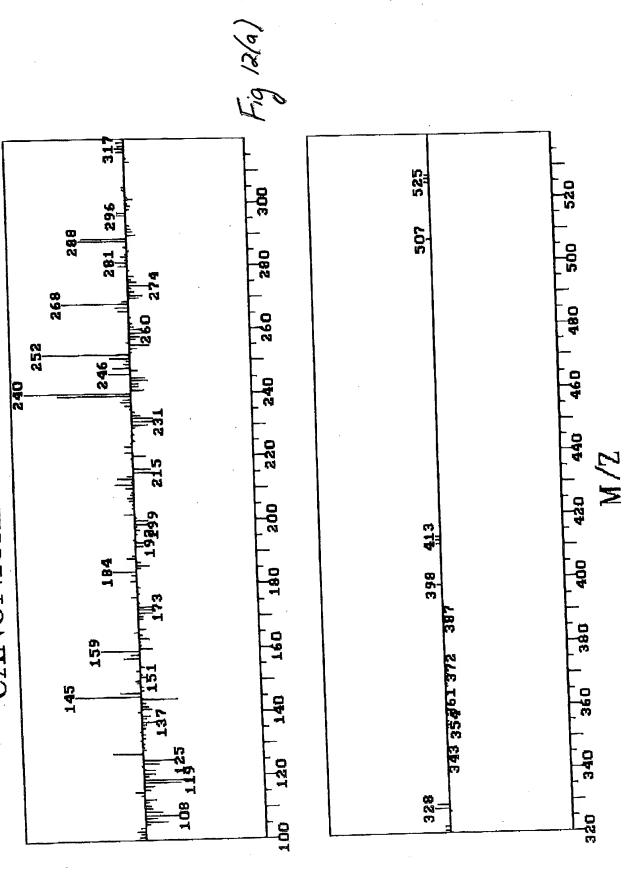


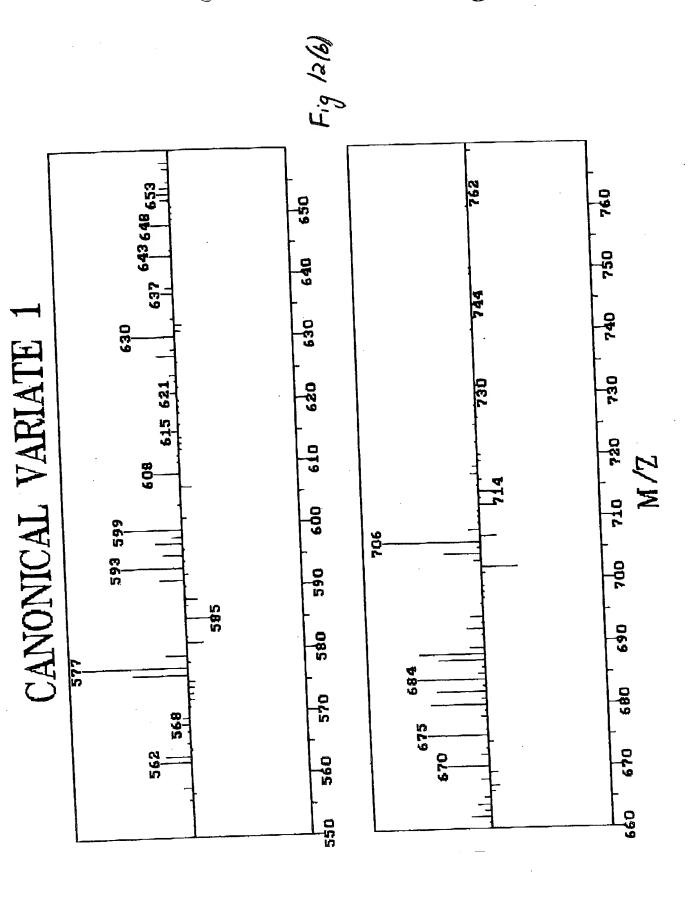






COMPONENT 1





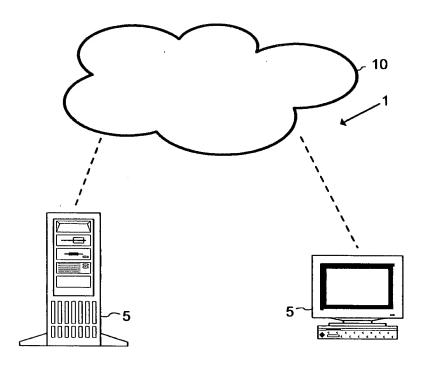


Fig 13

